

d hist

(FILE 'HOME' ENTERED AT 14:33:20 ON 07 JAN 2005)

FILE 'REGISTRY' ENTERED AT 14:33:36 ON 07 JAN 2005

L1 STRUCTURE UPLOADED
L2 2 S L1
L3 26 S L1 FULL

FILE 'CAPLUS, CAOLD, USPATFULL' ENTERED AT 14:34:38 ON 07 JAN 2005

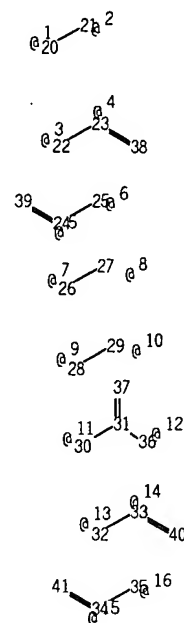
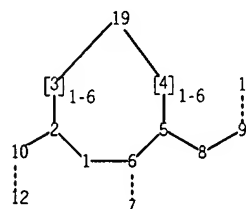
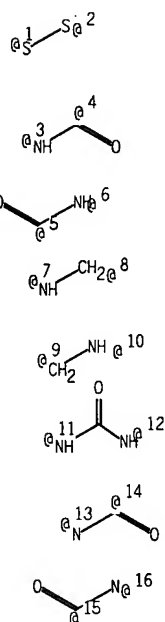
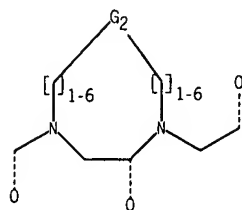
L4 2 S L3

=> d l4 1-2 cbib

L4 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2005 ACS on STN
2003:570952 Document No. 139:117689 Peptidyl heterocyclic compounds,
combinatorial libraries, and methods of selecting drug leads. Gilon,
Chaim (Yissum Research Development Company of the Hebrew University of
Jerusalem, Israel). PCT Int. Appl. WO 2003059876 A2 20030724, 71 pp.
DESIGNATED STATES: W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ,
CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE,
GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS,
LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO,
RU, SC, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC,
VN, YU, ZA, ZM, ZW; RW: AT, BE, BF, BJ, CF, CG, CH, CI, CM, CY, DE, DK,
ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD,
TG, TR. (English). CODEN: PIXXD2. APPLICATION: WO 2003-IL8 20030102.
PRIORITY: US 2002-34212 20020103.

L4 ANSWER 2 OF 2 USPATFULL on STN
2003:207899 Heterocyclic compounds, method of developing new drug leads and
combinatorial libraries used in such method.
Gilon, Chaim, Jerusalem, ISRAEL
YISSUM RESEARCH DEVELOPMENT COMPANY OF THE HEBREW UNIVERSITY OF JERUSALEM,
Jerusalem, ISRAEL (non-U.S. corporation)
US 2003144260 A1 20030731
APPLICATION: US 2002-34212 A1 20020103 (10)
DOCUMENT TYPE: Utility; APPLICATION.
CAS INDEXING IS AVAILABLE FOR THIS PATENT.

Search for claim 7



chain nodes :

7 8 9 10 11 12 20 21 22 23 24 25 26 27 28 29 30 31 32
33 34 35 36 37 38 39 40 41

ring nodes :

1 2 3 4 5 6 19

chain bonds :

2-10 5-8 6-7 8-9 9-11 10-12 20-21 22-23 23-38 24-25 24-39
26-27 28-29 30-31 31-36 31-37 32-33 33-40 34-35 34-41

ring bonds :

1-2 1-6 2-3 3-19 4-5 4-19 5-6

exact/norm bonds :

1-2 1-6 2-3 2-10 3-19 4-5 4-19 5-6 5-8 6-7 8-9 9-11 10-12
20-21 22-23 23-38 24-25 24-39 26-27 28-29 30-31 31-36 31-37
32-33 33-40 34-35 34-41

G1:C,O,S,N

G2:[*1-*2],[*3-*4],[*5-*6],[*7-*8],[*9-*10],[*11-*12],[*13-*14],[*15-*16]

Hydrogen count :

1:>= minimum 1 8:>= minimum 1

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 9:CLASS
10:CLASS 11:CLASS 12:CLASS 19:CLASS 20:CLASS 21:CLASS 22:CLASS
23:CLASS 24:CLASS 25:CLASS 26:CLASS 27:CLASS 28:CLASS 29:CLASS
30:CLASS 31:CLASS 32:CLASS 33:CLASS 34:CLASS 35:CLASS 36:CLASS
37:CLASS 38:CLASS 39:CLASS 40:CLASS 41:CLASS